## **Presently Pending Claims**

Language to be added has been <u>bolded and underlined</u>, while language to be deleted has been <del>bolded and striken-through</del>.

1-20. (Canceled)

21. (Currently amended) A compound represented by the structural formula I

$$R^{1}$$
 $R^{2}$ 
 $R^{9}$ 
 $R^{10}$ 
 $R^{10}$ 
 $R^{10}$ 
 $R^{3}$ 
 $R^{3}$ 

or a pharmaceutically acceptable salt or solvate thereof; wherein:

R<sup>1</sup> is

§—M—R<sup>4</sup> ;

R<sup>2</sup> is arylalkyl;

R<sup>3</sup> is selected from the group consisting of 6-membered heteroaryl, and 6-membered heteroaryl-N-oxide, wherein said 6-membered heteroaryl or heteroaryl-N-oxide is pyrimidine or pyrimidine-N-oxide respectively, each of which is optionally substituted with 1-4 substituents which can be the same or different and are independently selected from the group consisting of R<sup>11</sup>, R<sup>12</sup>, R<sup>13</sup>, R<sup>14</sup> and R<sup>15</sup>;

 $R^4$  is 1-3 substituents selected from the group consisting of H, halo, alkyl, haloalkyl, alkoxy, cycloalkyl, amide,  $CF_3$ ,  $OCF_3$ , aryl, heteroaryl,  $-XR^7$ , -CN,  $-CO_2H$ ,  $-CO_2R^{22}$ ,  $R^8$ -aryl( $C_1$ - $C_6$ )alkyl-,  $R^8$ -heteroaryl( $C_1$ - $C_6$ )alkyl-,  $-C(O)NR^{21}R^{22}$ ,  $-C(O)NH_2$ , wherein  $R^4$  can be the same or different and is independently selected when there is more than one  $R^4$  present;

R<sup>7</sup> is selected from the group consisting of aryl, substituted aryl, heteroaryl, alkyl, haloalkyl and cycloalkyl;

R<sup>8</sup> is 1, 2 or 3 substituents selected from the group consisting of H, halo, (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, -CF<sub>3</sub>, -OCF<sub>3</sub>, CH<sub>3</sub>C(O)-, -CN, CH<sub>3</sub>SO<sub>2</sub>-,

CF<sub>3</sub>SO<sub>2</sub>- and –NH<sub>2</sub>, wherein R<sup>8</sup> can be the same or different and is independently selected when there are more than one R<sup>8</sup> present;

 $R^9$ ,  $R^{10}$  and B can be the same or different and are each independently selected from the group consisting of hydrogen,  $(C_1-C_6)$ alkyl, and  $-(C_1-C_6)$ haloalkyl;

 $R^{11}$  and  $R^{12}$  can be the same or different and are each independently selected from the group consisting of ( $C_1$ - $C_6$ )alkyl, -( $C_1$ - $C_6$ )haloalkyl, halogen, -NR<sup>19</sup>R<sup>20</sup>, -OH, CF<sub>3</sub>, -OCH<sub>3</sub>, -O-acyl, and -OCF<sub>3</sub>;

 $R^{13}$  is selected from the group consisting of hydrogen,  $R^{11}$ , H, phenyl, -NO2, -CN, -CH2F, -CHF2, -CHO, -CH=NOR $^{19}$ , pyridyl-N-oxide, pyrimidinyl, pyrazinyl,  $N(R^{20})CONR^{20}R^{21}$ , -NHCONH(chloro-(C1-C6)alkyl), -NHCONH((C3-C10)-cycloalkyl(C1-C6)alkyl), -NHCO(C1-C6)alkyl, -NHCOCF3, -NHCOCF3, -NHSO2N((C1-C6)alkyl)2, -NHSO2(C1-C6)alkyl, -N(SO2CF3)2, -NHCO2(C1-C6)alkyl, (C3-C10)cycloalkyl, -SR $^{22}$ , -SOR $^{22}$ , -SO2R $^{22}$ , -SO2NH(C1-C6 alkyl), -OSO2(C1-C6)alkyl, -OSO2CF3, hydroxy(C1-C6)alkyl, -CONR $^{19}R^{20}$ , -CON(CH2CH2-O-CH3)2, -OCONH(C1-C6)alkyl, -CO2R $^{19}$ , -Si(CH3)3 and -B(OC(CH3)2)2;

 $R^{14}$  is selected from the group consisting of (C<sub>1</sub>-C<sub>6</sub>)alkyl, -(C<sub>1</sub>-C<sub>6</sub>)haloalkyl -NH<sub>2</sub> and  $R^{15}$ -phenyl;

 $R^{15}$  is 1-3 substituents selected from the group consisting of hydrogen, (C<sub>1</sub>-C<sub>6</sub>)alkyl, -(C<sub>1</sub>-C<sub>6</sub>)haloalkyl, -CF<sub>3</sub>, -CO<sub>2</sub>R<sup>20</sup>, -CN, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and halogen; wherein  $R^{15}$  can be the same or different and is independently selected when there are more than one  $R^{15}$  present;

 $R^{19}$ ,  $R^{20}$  and  $R^{21}$  can each be the same or different and are each independently selected from the group consisting of H,  $(C_1-C_6)$ alkyl and  $(C_3-C_6)$ cycloalkyl;

 $R^{22}$  is selected from the group consisting of  $(C_1-C_6)$ alkyl,  $-(C_1-C_6)$ haloalkyl,  $(C_2-C_6)$ hydroxyalkyl,  $(C_2-C_6)$ alkylene,  $(C_3-C_6)$ cycloalkyl, aryl and aryl $(C_1-C_6)$ alkyl-;

A is selected from the group consisting of H,  $(C_1-C_6)$ alkyl, and  $(C_2-C_6)$  alkenyl. M is aryl optionally substituted with  $R^4$ ; and

X is selected from the group consisting of  $\text{CH}_{2,}$  SO $_2,$  SO, S, and O, with the following proviso:

when  $R^1$  is phenyl, or naphthyl,  $R^2$  cannot be H, -(C<sub>1</sub>-C<sub>6</sub>)alkyl or -C(O)-(C<sub>1</sub>-C<sub>6</sub>)alkyl.

22. (Currently amended) A compound having the structural formula I according to claim 21 or a pharmaceutically acceptable salt **or solvate** thereof, wherein R<sup>9</sup>, R<sup>10</sup> and B are H, A is -CH<sub>3</sub>, and R<sup>1</sup>, R<sup>2</sup> and R<sup>3</sup> are as defined in the following table:

#	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>
8	Br		N
9	T'i		N N
10	CF <sub>3</sub>		N N
11	CF <sub>3</sub> O		N N
12	CI		N
14	F		N N
17	Br		N N N
18	CI		N
19	Br	F	N N N
20	CF <sub>3</sub>		Z Z Z

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·		Attorney Docket No
21		N
22		N N
23		N N
24	OCF <sub>3</sub>	N N
25	F F	N N N
26	CI	N N
27	F CI	N N
28	Br ← CH <sub>3</sub>	N N
29	MeO CI	N N
31	F Y	N 2
32	CI	N N N
33	OMe	N

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		 Attorney Docket No
34	F <sub>3</sub> C Y Y Y Y Y Y Y Y Y Y Y Y Y Y Y Y Y Y Y	N
35	CI CF <sub>3</sub>	N N N N N N N N N N N N N N N N N N N
36	F Yu	N N N
37	MeO	N N
38	Br CF <sub>3</sub>	N N
39	CH <sub>3</sub>	N N
40	EtO	N N
41	Et	N N
42	F	N N
43	PhO	N
44	CN	N N
50	MeSO <sub>2</sub>	N

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		Attorney Docket N
59	HO <sub>2</sub> C	N N N
60	HO <sub>2</sub> C	N N
61	H <sub>2</sub> N	N
63	O ZZ	Z Z Z
65	NH NH	N N N N N N N N N N N N N N N N N N N
66	O N H	N N
67	HO NH H	N N
68	A N H	N N N N N N N N N N N N N N N N N N N

23. (Currently amended) A compound according to claim 22 or a pharmaceutically acceptable salt thereof, wherein R<sup>1</sup>, R<sup>2</sup> and R<sup>3</sup> each represent:

#	R	R <sup>2</sup>	R <sup>3</sup>
10	CF <sub>3</sub>		N N
- Annual	CF <sub>3</sub> O		N N
12	CI		N

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		 Attorney Docket
14	F	N N
17	Br Br	N
28	Br CH <sub>3</sub>	N N
29	MeO CI	N N N N N N N N N N N N N N N N N N N
31	F	N N N N N N N N N N N N N N N N N N N
36	F F	N N
37	MeO	N N
39	CH <sub>3</sub>	N
40	EtO	, , , , , , , , , , , , , , , , , , ,
50	MeSO <sub>2</sub>	N N N N N N N N N N N N N N N N N N N
61	H <sub>2</sub> N	N N
68	ZH ZH	N N

24. (Currently amended) A compound according to claim 23 represented by the structural formulae:

and

or a pharmaceutically acceptable salt or solvate thereof.

- 25. (Currently amended) A pharmaceutical composition comprising one or more compounds of claim 21 or a pharmaceutically acceptable salt thereof.
- 26. (Currently amended) A pharmaceutical composition comprising one or more compounds of claim 24 or a pharmaceutically acceptable salt thereof.
- 27. (Previously presented) The pharmaceutical composition according to claim 25 further comprising one or more pharmaceutically acceptable carriers.
- 28. (Previously presented) The pharmaceutical composition according to claim 26 further comprising one or more pharmaceutically acceptable carriers.
- 29. (Currently amended) The pharmaceutical composition according to claim25, wherein said pharmaceutical composition contains a therapeutically effective

- 28. (Previously presented) The pharmaceutical composition according to claim 26 further comprising one or more pharmaceutically acceptable carriers.
- 29. (Currently amended) The pharmaceutical composition according to claim 25, wherein said pharmaceutical composition contains a therapeutically effective amount of said one or more compounds or a pharmaceutically acceptable salt **er solvate** thereof.
- 30. (Currently amended) The pharmaceutical composition according to claim 26, wherein said pharmaceutical composition contains a therapeutically effective amount of said one or more compounds or a pharmaceutically acceptable salt **or solvate** thereof.

31-40. (Canceled)